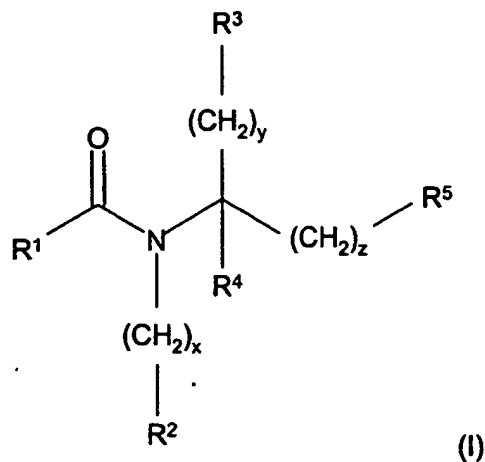


IN THE CLAIMS

1. (canceled).

2. (Previously presented) A compound of formula (I)



wherein:

R¹ is selected from:

- a) phenyl, which is optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, CF₃, halo, CN, NR⁷R⁸, SO₂R⁶ and OC₁-C₆ alkyl, and
- b) Aromatic Heterocycle, wherein said Aromatic Heterocycle is selected from pyridyl, pyrazinyl, pyrimidinyl, quinolinyl, quinoxalinyl, isoxazolyl and pyrazolyl, each aromatic heterocycle optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, SR⁶, SO₂R⁶, NH₂, CF₃, halo, OH, OC₁-C₆ alkyl, NR⁷R⁸ wherein R⁸ may be optionally substituted by NH₂, phenyl or Heterocycle, and OPh wherein Ph may be optionally substituted by 1-3 groups each independently selected from halo and C₁-C₆ alkyl;

R² is selected from:

- a) phenyl, which is optionally substituted by methyl, fluoro, chloro, methoxy, CF₃ or SO₂CH₃,
- b) pyrazolyl, which is optionally substituted by methyl, and
- c) C(O)N(CH₃)₂;

R³ is selected from:

- a) phenyl, said phenyl being optionally fused to Heterocycle and said phenyl or said fused phenyl being optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, halo, CN and OC₁-C₆ alkyl,
- b) R⁶,
- c) cyclopropyl, cyclobutyl, cyclopentyl or cyclohexyl, which is optionally substituted by C₁-C₆ alkyl; and
- d) Aromatic Heterocycle, wherein said Aromatic Heterocycle may be defined as a 5-6 membered aromatic heterocycle containing 1 or 2 nitrogen atoms, said ring optionally fused with a phenyl or a 3-8 membered cycloalkyl group.

R⁴ is H;

R⁵ is CONH₂;

R⁶ is methyl;

R⁷ is hydrogen or C₁-C₆ alkyl;

R⁸ is C₁-C₆ alkyl;

or NR⁷R⁸ forms a monocyclic saturated ring system containing between 3 and 7 ring atoms;

x is 1;

y is 0; and

z is 0 or 1

wherein:

Aromatic Heterocycle may be defined as a 5-6 membered aromatic heterocycle containing 1-4 heteroatoms each independently selected from N, O and S, said ring optionally fused with a phenyl or a 3-8 membered cycloalkyl group;

Heterocycle is a 5-8 membered saturated or partially saturated ring containing 1-3 heteroatoms each independently selected from N, O and S, said ring optionally fused with phenyl;

a tautomer thereof or a pharmaceutically acceptable salt, solvate or polymorph of said compound or tautomer.

2. (Currently amended) A compound according to claim 1 wherein R¹ is selected from:

- a) phenyl, which is optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, CF₃, halo, CN, NR⁷R⁸, SO₂R⁶ and OC₁-C₆ alkyl, and
- b) Aromatic Heterocycle, wherein said Aromatic Heterocycle is selected from:
 - i) pyridyl, which is optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, SO₂R⁶, NH₂, CF₃, CN, halo, OH, OC₁-C₆ alkyl, NR⁷R⁸ wherein R⁸ may be optionally substituted by NH₂, phenyl or Heterocycle, and OPh wherein Ph may be optionally substituted by 1-3 groups each independently selected from halo and C₁-C₆ alkyl;
 - ii) pyrimidinyl, which is optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, SO₂R⁶, NH₂, CF₃, CN, halo, OH, OC₁-C₆ alkyl, NR⁷R⁸ wherein R⁸ may be optionally substituted by NH₂, phenyl or Heterocycle, and OPh wherein Ph may be optionally substituted by 1-3 groups each independently selected from halo and C₁-C₆ alkyl;
 - iii) pyrazinyl, which is optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, NH₂, SR⁶ and halo;
 - iv) quinolinyl;
 - v) quinoxalinyl, which is optionally substituted by OH;
 - vi) isoxazolyl, which is optionally substituted by 1-3 groups each independently selected from: C₁-C₆ alkyl; and
 - vii) pyrazole;

~~R² is selected from:~~

- ~~a) phenyl, which is optionally substituted by methyl, halo, methoxy, CF₃ or SO₂CH₃;~~
- ~~b) cyclopropyl or 1- or 2-indanyl;~~
- ~~c) pyrazolyl, which is optionally substituted by methyl;~~
- ~~d) C(O)N(CH₃)₂, and~~
- ~~e) piperidinyl substituted by C(O)R⁶.~~

R³ is selected from:

- a) phenyl, said phenyl being optionally fused to 1,4-dioxan and said phenyl or said fused phenyl being optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, halo, CN and OC₁-C₆ alkyl;
- b) R⁶,
- c) cyclopropyl, which is optionally substituted by C₁-C₆ alkyl; and
- d) Aromatic Heterocycle, wherein said Aromatic Heterocycle is selected from pyrazolyl or pyridyl, both optionally substituted by C₁-C₆ alkyl;

~~R⁵ is CONH₂ or CH₃;~~ and

z is 0.

³
~~2~~ ⁴ (Currently amended) A compound according to any one of claims 1 to ~~2~~ or ~~3~~ wherein R¹ is phenyl, 2- or 3-pyridyl or 2,4-pyrimidinyl, said moieties being optionally substituted by 1-3 groups each independently selected from C₁-C₆ alkyl, halo, OC₁-C₆ alkyl, CN, SO₂R⁶, NHR₇, NHCH₂CH₂NH₂ and CF₃;

⁴
~~5~~ (original) A compound according to claim ~~4~~ ³ wherein R¹ is phenyl, 2- or 3-pyridyl or 2,4-pyrimidinyl, said moieties being optionally substituted by 1-3 groups each independently selected from methyl, fluoro, chloro, methoxy, ethoxy, n-propoxy, CN, SO₂CH₃, NH₂, NHCH₃, NHCH₂CH₂NH₂, and CF₃.

6. (canceled)

⁵
~~7~~ (previously presented) A compound according to claim ~~5~~ ⁴ wherein R² is phenyl, *para*-fluorophenyl, *para*-chlorophenyl, *para*-methylphenyl, 2,5-dimethylphenyl, *o*-methylphenyl and *para*-methoxyphenyl.

⁵
~~6~~ (previously presented) A compound according to claim ~~7~~ ⁵ wherein R³ is selected from:

- a) phenyl, said phenyl being optionally fused to 1,4-dioxan and said phenyl or said fused phenyl being optionally substituted by 1-2 groups each independently selected from methyl, methoxy, ethoxy, fluoro, chloro and CN;

- b) isopropyl;
- c) cyclopropyl; and
- d) pyrazolyl and pyridyl, both optionally substituted by methyl.

7 ~~9~~ (original) A compound according to claim ~~8~~⁶ wherein R³ is 3-methoxyphenyl or 1,4-benzodioxanyl.

10. (Cancelled).

8 ~~11~~ (Previously presented) A compound according to claim ~~2~~¹ selected from:

2-Amino-N-[2-amino-1-(2-methylphenyl)-2-oxoethyl]-N-(4-chlorobenzyl)nicotinamide,

N-[2-Amino-1-(3-methoxyphenyl)-2-oxoethyl]-4-cyano-N-(4-methylbenzyl)benzamide,

N-[3-Amino-1-(3-methoxyphenyl)-3-oxopropyl]-4-methyl-N-(4-methylbenzyl)nicotinamide,

2-Amino-N-[(1S)-3-amino-3-oxo-1-phenylpropyl]-N-(4-methylbenzyl)nicotinamide,

5-Chloro-2-methylthio-N-[2-amino-1-{1,4-benzodioxan-6-yl}-2-oxoethyl]-N-(4-methylbenzyl)pyrimidine-4-carboxamide,

5-Chloro-2-amino-N-[2-amino-1-{1,4-benzodioxan-6-yl}-2-oxoethyl]-N-(4-methylbenzyl)pyrimidine-4-carboxamide, and

2-Amino-N-[carbamoyl-(2,3-dihydro-benzo[1,4]dioxin-6-yl)-methyl]-4,6-dimethyl-N-(4-methyl-benzyl)-nicotinamide;

and tautomers thereof and pharmaceutically acceptable salts, solvates and polymorphs of said compound or tautomer.

9 ~~12~~ (Previously presented) A pharmaceutical composition comprising a compound of claim ~~2~~¹, or pharmaceutically acceptable salts, solvates or polymorphs thereof, and a pharmaceutically acceptable diluent or carrier.

~~10~~ ~~13~~ (previously canceled)

~~10~~ ~~14~~ (Previously presented) A method of treatment of a disorder or condition where inhibition of Oxytocin is known, or can be shown, to produce a beneficial